

(I)

Y is oxygen, NR_{4a}, sulfur, sulfonyl, sulfinyl, C(O), C(=NR_{4b}), C(=CR_{6a}R_{6b}) or a C₁-C₄alkylene or C₂-C₄alkenylene chain, which may be interrupted by oxygen, NR_{5a}, sulfur, sulfonyl, sulfinyl, C(O) or C(=NR_{5b}) and/or mono- or poly-substituted by R₆;

A_2 is nitrogen or CR_8 .

R₁, R₂, R₆, R₇ and R₈ are each independently of the others hydrogen, hydroxy, mercapto, NO₂, cyano, halogen, formyl, oxyiminomethylene, C₁-C₆alkoxyiminomethylene, C₁-C₆alkyl, C₁-C₆haloalkyl, C₂-C₆alkenyl, C₂-C₆haloalkenyl, C₂-C₆alkynyl, C₂-C₆haloalkynyl, C₁-C₆alkoxy, C₁-C₆haloalkoxy, C₃-C₆alkenyloxy, C₃-C₆alkynyloxy, C₃-C₆oxacycloalkyl, C₃-C₆thiacycloalkyl, C₃-C₆dioxacycloalkyl, C₃-C₆dithiacycloalkyl, C₃-C₆oxathiacycloalkyl, C₁-C₆alkoxycarbonyl, C₁-C₆alkylcarbonyl, C₁-C₆alkoxycarbonyloxy, C₁-C₆alkylcarbonyloxy, C₁-C₆alkylthio, C₁-C₆alkylsulfonyl, C₁-C₆alkylsulfinyl, NR₉R₁₀, C₃-C₆cycloalkyl, tri(C₁-C₆alkyl)silyl, di(C₁-C₆alkyl)phenylsilyl, tri(C₁-C₆alkyl)silyloxy, di(C₁-C₆alkyl)phenylsilyloxy or Ar₁;

or R₁, R₂, R₆, R₇, R₈ are each independently of the others a C₁-C₆alkyl, C₂-C₆alkenyl, C₂-C₆alkynyl or C₃-C₆cycloalkyl group, which may be interrupted by oxygen, sulfur, sulfonyl, sulfinyl, -NR₁₁- or -C(O)- and/or mono-, di- or tri-substituted by hydroxy, mercapto, NO₂, cyano, halogen, formyl, C₁-C₆alkoxy, C₃-C₆alkenyloxy, C₃-C₆alkynyloxy, C₁-C₆haloalkoxy, C₁-C₂alkoxy-C₁-C₂alkoxy, C₁-C₄alkoxycarbonyloxy, C₁-C₄alkylcarbonyloxy, C₁-C₄alkoxy-carbonyl, C₁-C₄alkylcarbonyl, C₁-C₆alkylthio, C₁-C₆alkylsulfinyl, C₁-C₆alkylsulfonyl, NR₁₂R₁₃, C₁-C₆alkyl, C₂-C₆alkenyl, C₂-C₆alkynyl, C₃-C₆cycloalkyl, tri(C₁-C₆alkyl)silyl, tri(C₁-C₆alkyl)-silyloxy or Ar₂;

or two substituents R₆ at the same carbon atom together form a –CH₂O– or a C₂–C₅alkylene chain, which may be interrupted once or twice by oxygen, sulfur, sulfinyl or sulfonyl and/or mono- or poly-substituted by R_{6c}, with the proviso that two hetero atoms may not be located next to one another;

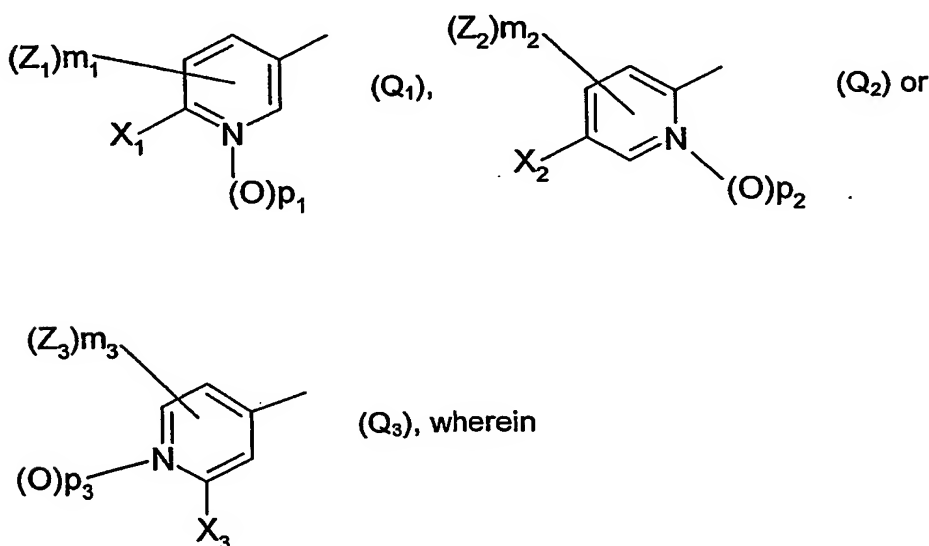
or two substituents R_6 at different carbon atoms together form an oxygen bridge or a C_1 - C_4 alkylene chain, which may in turn be substituted by R_{6ci} ;

or R_7 and R_8 together form a $-\text{CH}_2\text{CH}=\text{CH}-$, $-\text{OCH}=\text{CH}-$ or $-\text{CH}=\text{CH}-\text{CH}=\text{CH}-$ bridge or a C_3 - C_4 alkylene chain, which may be interrupted by oxygen or $-\text{S}(\text{O})_{n1}-$ and/or mono- or poly-substituted by R_{6d} ;

R_3 is hydroxy, halogen, mercapto, C_1 - C_8 alkylthio, C_1 - C_8 alkylsulfinyl, C_1 - C_8 alkylsulfonyl, C_1 - C_8 haloalkylthio, C_1 - C_8 haloalkylsulfinyl, C_1 - C_8 haloalkylsulfonyl, C_1 - C_4 alkoxy- C_1 - C_4 alkylthio, C_1 - C_4 alkoxy- C_1 - C_4 alkylsulfinyl, C_1 - C_4 alkoxy- C_1 - C_4 alkylsulfonyl, C_3 - C_8 alkenylthio, C_3 - C_8 alkynylthio, C_1 - C_4 alkylthio- C_1 - C_4 alkylthio, C_3 - C_4 alkenylthio- C_1 - C_4 alkylthio, C_1 - C_4 alkoxy-carbonyl- C_1 - C_4 alkylthio, C_1 - C_4 alkoxycarbonyl- C_1 - C_4 alkylsulfinyl, C_1 - C_4 alkoxycarbonyl- C_1 - C_4 alkylsulfonyl, C_3 - C_8 cycloalkylthio, C_3 - C_8 cycloalkylsulfinyl, C_3 - C_8 cycloalkylsulfonyl, phenyl- C_1 - C_4 alkylthio, phenyl- C_1 - C_4 alkylsulfinyl, phenyl- C_1 - C_4 alkylsulfonyl, $\text{S}(\text{O})_{n1}-\text{Ar}_3$, phenylthio, phenylsulfinyl, phenylsulfonyl, it being possible for the phenyl-containing groups to be substituted by one or more C_1 - C_3 alkyl, C_1 - C_3 haloalkyl, C_1 - C_3 alkoxy, C_1 - C_3 haloalkoxy, C_1 - C_4 alkoxycarbonyl, halogen, cyano, hydroxy or nitro groups;

or R_3 is O^+M^+ , wherein M^+ is an alkali metal cation or an ammonium cation;

Q is a radical



p_1 , p_2 and p_3 are 0 or 1;

m_1 , m_2 and m_3 are 1, 2 or 3;

X_1 , X_2 and X_3 are hydroxy, halogen, C_1 - C_6 alkyl, C_1 - C_6 haloalkyl, C_2 - C_6 alkenyl, C_2 - C_6 haloalkenyl, C_2 - C_6 alkynyl, C_2 - C_6 haloalkynyl, C_1 - C_6 alkoxy, C_1 - C_6 haloalkoxy, C_1 - C_6 alkylthio, C_1 - C_6 alkylsulfinyl, C_1 - C_6 alkylsulfonyl, C_1 - C_6 haloalkylthio, C_1 - C_6 haloalkylsulfinyl or C_1 - C_6 haloalkylsulfonyl;

Z_1 , Z_2 and Z_3 are C_1 - C_6 alkyl which is substituted by the following substituents: C_3 - C_4 cycloalkyl or C_3 - C_4 cycloalkyl substituted by halogen, C_1 - C_6 alkyl, C_1 - C_3 alkoxy or C_1 - C_3 haloalkoxy-

C_1-C_3 alkyl; oxiranyl or oxiranyl substituted by C_1-C_6 alkyl or C_1-C_3 alkoxy- C_1-C_3 alkyl; 3-oxetanyl or 3-oxetanyl substituted by C_1-C_6 alkyl, C_1-C_3 alkoxy or C_1-C_3 alkoxy- C_1-C_3 alkyl; 3-oxetanyloxy or 3-oxetanyloxy substituted by C_1-C_6 alkyl, C_1-C_3 alkoxy or C_1-C_3 alkoxy- C_1-C_3 alkyl; C_3-C_6 cycloalkyloxy or C_3-C_4 cycloalkyloxy substituted by halogen, C_1-C_6 alkyl, C_1-C_3 alkoxy or C_1-C_3 alkoxy- C_1-C_3 alkyl; C_1-C_6 haloalkoxy; C_1-C_6 alkylsulfonyloxy; C_1-C_6 haloalkylsulfonyloxy; phenylsulfonyloxy; benzylsulfonyloxy; benzoyloxy; phenoxy; phenylthio; phenylsulfinyl; phenylsulfonyl; Ar_{10} ; OAr_{12} ; $tri(C_1-C_6alkyl)silyl$ or $tri(C_1-C_6alkyl)silyloxy$, it being possible for the phenyl-containing groups to be mono- or poly-substituted by C_1-C_3 alkyl, C_1-C_3 haloalkyl, C_1-C_3 alkoxy, C_1-C_3 haloalkoxy, halogen, cyano, hydroxy or nitro; or Z_1 , Z_2 and Z_3 are 3-oxetanyl; 3-oxetanyl substituted by C_1-C_3 alkoxy, C_1-C_3 alkoxy- C_1-C_3 alkyl or C_1-C_6 alkyl; C_3-C_6 cycloalkyl substituted by halogen, C_1-C_3 alkyl or C_1-C_3 alkoxy- C_1-C_3 alkyl; $tri(C_1-C_6alkyl)silyl$; $tri(C_1-C_6alkyl)silyloxy$ or $CH=P(phenyl)_3$; or Z_1 , Z_2 and Z_3 are a C_1-C_6 alkyl, C_2-C_6 alkenyl or C_2-C_6 alkynyl group, which is interrupted by oxygen, $-O(CO)-$, $-(CO)O-$, $-O(CO)O-$, $-N(R_{14})O-$, $-ONR_{15}-$, sulfur, sulfinyl, sulfonyl, $-SO_2NR_{16}-$, $-NR_{17}SO_2-$ or $-NR_{18}-$ and is mono- or poly-substituted by L_1 ; it also being possible for L_1 to be bonded at the terminal carbon atom of the C_1-C_6 alkyl, C_2-C_6 alkenyl or C_2-C_6 alkynyl group; or Z_1 , Z_2 and Z_3 are hydrogen, hydroxy, mercapto, NO_2 , cyano, halogen, formyl, C_1-C_6 alkyl, C_1-C_6 haloalkyl, C_2-C_6 alkenyl, C_2-C_6 haloalkenyl, C_2-C_6 alkynyl, C_2-C_6 haloalkynyl, C_1-C_6 alkoxy, C_1-C_6 haloalkoxy, C_1-C_6 alkoxycarbonyl, C_1-C_6 alkylcarbonyl, C_1-C_6 alkylthio, C_1-C_6 alkylsulfonyl, C_1-C_6 alkylsulfinyl, $NR_{22}R_{23}$, phenyl which may be mono- or poly-substituted by C_1-C_3 alkyl, C_1-C_3 haloalkyl, C_1-C_3 alkoxy, C_1-C_3 haloalkoxy, halogen, cyano, hydroxy or nitro, C_3-C_6 cycloalkyl, C_5-C_6 cycloalkyl substituted by C_1-C_3 alkoxy, C_1-C_3 alkoxy- C_1-C_3 alkyl or C_1-C_6 alkyl, or Ar_5 , $O-Ar_6$, $N(R_{24})Ar_7$ or $S(O)_nAr_8$; L_1 is hydrogen, halogen, hydroxy, amino, formyl, nitro, cyano, mercapto, carbamoyl, $P(O)(OC_1-C_6alkyl)_2$, C_1-C_6 alkoxy, C_1-C_6 haloalkoxy, C_1-C_6 alkoxycarbonyl, C_2-C_6 alkenyl, C_2-C_6 haloalkenyl, C_2-C_6 alkynyl, C_2-C_6 haloalkynyl, C_3-C_6 cycloalkyl, halo-substituted C_3-C_6 cycloalkyl, C_3-C_6 alkenyloxy, C_3-C_6 alkynyloxy, C_3-C_6 haloalkenyloxy, cyano- C_1-C_6 alkoxy, C_1-C_6 alkoxy- C_1-C_6 alkoxy, C_1-C_6 alkylthio- C_1-C_6 alkoxy, C_1-C_6 alkylsulfinyl- C_1-C_6 alkoxy, C_1-C_6 alkylsulfonyl- C_1-C_6 alkoxy, C_1-C_6 alkoxycarbonyl- C_1-C_6 alkoxy, C_1-C_6 alkylcarbonyloxy- C_1-C_6 alkylcarbonyl, C_1-C_6 alkylthio, C_1-C_6 alkylsulfinyl, C_1-C_6 alkylsulfonyl, C_1-C_6 haloalkylthio, C_1-C_6 haloalkylsulfinyl, C_1-C_6 haloalkylsulfonyl or oxiranyl, which may in turn be substituted by C_1-C_6 alkyl, C_1-C_3 alkoxy or C_1-C_3 alkoxy- C_1-C_3 alkyl, or (3-oxetanyl)-oxy, which may in turn be substituted by C_1-C_6 alkyl, C_1-C_3 alkoxy or C_1-C_3 alkoxy- C_1-C_3 alkyl, or benzoyloxy, benzyloxy, benzylthio, benzylsulfinyl, benzylsulfonyl, C_1-C_6 alkylamino, $di(C_1-C_6alkyl)amino$, $R_{19}S(O)_2O-$, $R_{20}N(R_{21})SO_2-$, rhodano, phenyl, phenoxy, phenylthio, phenylsulfinyl, phenylsulfonyl, Ar_4 or

OAr₁₁, it being possible for the phenyl-containing groups in turn to be substituted by one or more C₁-C₃alkyl, C₁-C₃haloalkyl, C₁-C₃alkoxy, C₁-C₃haloalkoxy, halogen, cyano, hydroxy or nitro groups;

R_{4a} and R_{5a} are each independently of the other hydrogen, C₁-C₆alkyl, C₁-C₆haloalkyl, cyano, formyl, C₁-C₆alkylcarbonyl, C₁-C₆alkoxycarbonyl, carbamoyl, C₁-C₆alkylaminocarbonyl, di(C₁-C₆alkylamino)carbonyl, di(C₁-C₆alkylamino)sulfonyl, C₃-C₆cycloalkylcarbonyl, C₁-C₆alkylsulfonyl, phenylcarbonyl, phenylaminocarbonyl or phenylsulfonyl, it being possible for the phenyl groups to be mono- or poly-substituted by C₁-C₆alkyl, C₁-C₆haloalkyl, C₁-C₆alkoxy, C₁-C₆haloalkoxy, halogen, cyano, hydroxy or nitro;

R_{4b} and R_{5b} are each independently of the other hydroxy, C₁-C₆alkoxy, C₃-C₆alkenyloxy, C₃-C₆alkynyloxy or benzyloxy, it being possible for the benzyl group to be mono- or poly-substituted by C₁-C₆alkyl, C₁-C₆haloalkyl, C₁-C₆alkoxy, C₁-C₆haloalkoxy, halogen, cyano, hydroxy or nitro;

R₉, R₁₁, R₁₃, R₁₆, R₁₇, R₁₈, R₂₀, R₂₃ and R₂₄ are each independently of the others hydrogen, C₁-C₆alkyl, Ar₉, C₁-C₆haloalkyl, C₁-C₆alkylcarbonyl, C₁-C₆alkoxycarbonyl, C₁-C₆alkylsulfonyl, phenyl, it being possible for the phenyl group in turn to be mono- or poly-substituted by C₁-C₆alkyl, C₁-C₆haloalkyl, C₁-C₆alkoxy, C₁-C₆haloalkoxy, halogen, cyano, hydroxy or nitro; R_{6a} is hydrogen, C₁-C₆alkyl or C₁-C₆alkylcarbonyl; or together with R_{6b} is a C₂-C₅alkylene chain;

R_{6b}, R_{6d}, R₁₀, R₁₂ and R₂₂ are each independently of the others hydrogen or C₁-C₆alkyl;

R_{6c}, R₁₄, R₁₅, R₁₉ and R₂₁ are each independently of the others C₁-C₆alkyl or C₁-C₆haloalkyl;

Ar₁, Ar₂, Ar₃, Ar₄, Ar₅, Ar₆, Ar₇, Ar₈, Ar₉, Ar₁₀, Ar₁₁ and Ar₁₂ are each independently of the others a five- to ten-membered, monocyclic or fused bicyclic ring system, which may be aromatic, partially saturated or fully saturated and may contain from 1 to 4 hetero atoms selected from nitrogen, oxygen, sulfur, C(O) and C(=NR₂₅), and each ring system may contain not more than two oxygen atoms, not more than two sulfur atoms, not more than two C(O) groups and not more than one C(=NR₂₅) group, and each ring system may itself be mono- or poly-substituted by C₁-C₆alkyl, C₁-C₆haloalkyl, C₂-C₆alkenyl, C₂-C₆haloalkenyl, C₂-C₆alkynyl, C₂-C₆haloalkynyl, C₁-C₆alkoxy, C₁-C₆haloalkoxy, C₃-C₆alkenyloxy, C₃-C₆alkynyl-oxy, mercapto, amino, hydroxy, C₁-C₆alkylthio, C₁-C₆haloalkylthio, C₃-C₆alkenylthio, C₃-C₆haloalkenylthio, C₃-C₆alkynylthio, C₁-C₃alkoxy-C₁-C₃alkylthio, C₁-C₄alkylcarbonyl-C₁-C₃alkylthio, C₁-C₄alkoxycarbonyl-C₁-C₃alkylthio, cyano-C₁-C₃alkylthio, C₁-C₆alkylsulfinyl, C₁-C₆haloalkylsulfinyl, C₁-C₆alkylsulfonyl, C₁-C₆haloalkylsulfonyl, aminosulfonyl, C₁-C₂alkylamino-sulfonyl, N,N-di(C₁-C₂alkyl)aminosulfonyl, di(C₁-C₄alkyl)amino, halogen, cyano, nitro or phenyl, it being possible for the phenyl group in turn to be substituted by hydroxy, C₁-C₆alkylthio, C₁-C₆haloalkylthio, C₃-C₆alkenylthio, C₃-C₆haloalkenylthio, C₃-C₆alkynylthio,

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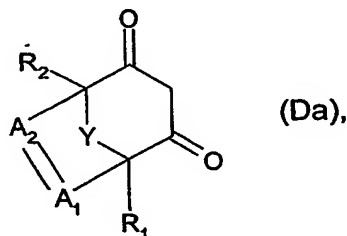
C₁-C₃alkoxy-C₁-C₃alkylthio, C₁-C₄alkylcarbonyl-C₁-C₃alkylthio, C₁-C₄alkoxycarbonyl-C₁-C₃alkylthio, cyano-C₁-C₃alkylthio, C₁-C₆alkylsulfinyl, C₁-C₆haloalkylsulfinyl, C₁-C₆alkylsulfonyl, C₁-C₆haloalkylsulfonyl, aminosulfonyl, C₁-C₂alkylaminosulfonyl, N,N-di(C₁-C₂alkyl)amino-sulfonyl, di(C₁-C₄alkyl)amino, halogen, cyano or nitro, and the substituents at the nitrogen atom in the heterocyclic ring being other than halogen, and two oxygen atoms not being located next to one another;

R₂₅ is hydrogen, hydroxy, C₁-C₆alkyl, C₁-C₆haloalkyl, C₁-C₆alkoxy, C₁-C₆alkylcarbonyl, C₁-C₆alkoxycarbonyl or C₁-C₆alkylsulfonyl; and

n₁ is 0, 1 or 2; and n₆ is 0, 1 or 2;

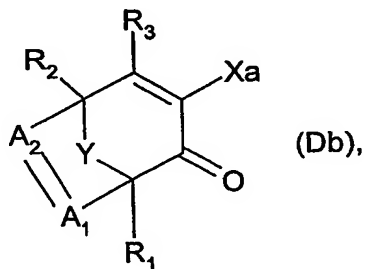
or an agronomically acceptable salt/isomer/enantiomer/tautomer of such a compound.

2. A compound of formula Da



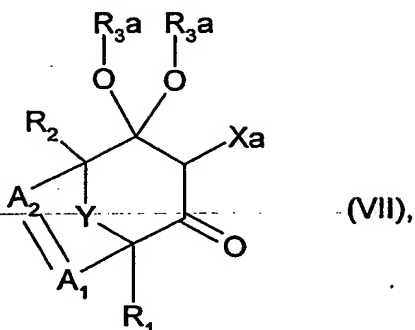
wherein Y, R₁, R₂, A₁ and A₂ are as defined for formula I in claim 1.

3. A compound of formula Db



wherein A₁, A₂, R₁, R₂ and Y are as defined for formula I in claim 1, Xa is hydrogen, chlorine or bromine and R₃ is hydroxy or C₁-C₆alkoxy, with the exception of the compounds 3-chloro-8-oxa-bicyclo[3.2.1]oct-6-ene-2,4-dione; 3-chloro-bicyclo[3.2.1]oct-6-ene-2,4-dione; 3-chloro-4-hydroxy-bicyclo[3.2.1]octa-3,6-dien-2-one; 3,4-dibromo-8-oxa-bicyclo[3.2.1]octa-3,6-dien-2-one; 3,4-dibromo-1,5-dimethyl-8-oxa-bicyclo[3.2.1]octa-3,6-dien-2-one; 3,4-dibromo-bicyclo[3.2.1]octa-3,6-dien-2-one; 3,4-dichloro-8-oxa-bicyclo[3.2.1]octa-3,6-dien-2-one; 3,4-dichloro-bicyclo[3.2.1]octa-3,6-dien-2-one and 7,8-dibromo-5,9-dihydro-5,9-methano-benzocyclohepten-6-one.

4. A compound of formula VII



wherein A_1 , A_2 , R_1 , R_2 , Y are as defined for formula I in claim 1, X_a is hydrogen, chlorine or bromine and R_{3a} is C_1 - C_6 alkyl or two R_{3a} together are $-CH_2CH_2-$.

5. A herbicidal and plant-growth-inhibiting composition, comprising a herbicidally effective amount of a compound of formula I according to claim 1 on an inert carrier.
6. A method of controlling undesired plant growth, which method comprises applying a compound of formula I according to claim 5, or a composition comprising such a compound, in a herbicidally effective amount to a plant or to the locus thereof.
7. A method of inhibiting plant growth, which method comprises applying a compound of formula I according to claim 5, or a composition comprising such a compound, in a herbicidally effective amount to a plant or to the locus thereof.